

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE030819\
 Data File : BE098844.D
 Acq On : 8 Mar 2019 22:17
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 09 05:07:39 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE030619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 06 15:33:24 2019
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	120	0.00
2	1,4-Dioxane	0.400	0.382	4.5	119	0.00
3	n-Nitrosodimethylamine	0.400	0.340	15.0	101	0.01
4 S	2-Fluorophenol	0.400	0.376	6.0	109	0.01
5 S	Phenol-d6	0.400	0.385	3.8	116	0.00
6	bis(2-Chloroethyl)ether	0.400	0.314	21.5	98	0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	125	0.00
8 S	Nitrobenzene-d5	0.400	0.358	10.5	125	0.00
9	Naphthalene	0.400	0.397	0.8	128	-0.01
10	Hexachlorobutadiene	0.400	0.391	2.3	126	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.396	1.0	127	0.00
12	2-Methylnaphthalene	0.400	0.388	3.0	125	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	132	0.00
14 S	2,4,6-Tribromophenol	0.400	0.390	2.5	130	-0.01
15 S	2-Fluorobiphenyl	0.400	0.383	4.3	128	-0.01
16	Acenaphthylene	0.400	0.382	4.5	133	0.00
17	Acenaphthene	0.400	0.389	2.8	129	-0.01
18	Fluorene	0.400	0.383	4.3	130	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	129	-0.01
20	4-Bromophenyl-phenylether	0.400	0.363	9.3	119	0.00
21	Hexachlorobenzene	0.400	0.368	8.0	125	0.00
22	Pentachlorophenol	0.400	0.419	-4.7	114	0.00
23	Phenanthrene	0.400	0.376	6.0	124	-0.01
24	Anthracene	0.400	0.352	12.0	120	0.00
25 SURR	Fluoranthene-d10	0.400	0.350	12.5	117	0.00
26	Fluoranthene	0.400	0.355	11.3	119	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	119	-0.01
28	Pyrene	0.400	0.398	0.5	119	0.00
29 S	Terphenyl-d14	0.400	0.380	5.0	114	0.00
30	Benzo(a)anthracene	0.400	0.355	11.3	108	0.00
31	Chrysene	0.400	0.393	1.8	120	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.337	15.8	110	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.391	2.3	119	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	125	0.00
35	Benzo(b)fluoranthene	0.400	0.381	4.8	123	0.00
36	Benzo(k)fluoranthene	0.400	0.379	5.3	124	-0.01
37 C	Benzo(a)pyrene	0.400	0.387	3.3	124	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.380	5.0	120	0.00
39	Benzo(g,h,i)perylene	0.400	0.381	4.8	121	0.00

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(#) = Out of Range	SPCC's out = 0 CCC's out = 0				